Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 13 SEP 2005 HIGHEST RN 863091-33-2 DICTIONARY FILE UPDATES: 13 SEP 2005 HIGHEST RN 863091-33-2

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

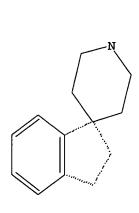
Structure search iteration limits have been increased. See HELP SLIMITS for details.

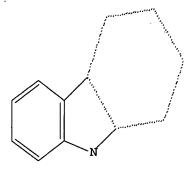
Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> Uploading C:\Program Files\Stnexp\Queries\10669606.str

L1 STRUCTURE UPLOADED

=> d l1 L1 HAS NO ANSWERS L1 STR





Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 13:41:54 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 535 TO ITERATE

100.0% PROCESSED 535 ITERATIONS 3 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 9313 TO 12087
PROJECTED ANSWERS: 3 TO 163

L2 3 SEA SSS SAM L1

=> s l1 ful

FULL SEARCH INITIATED 13:41:58 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 10612 TO ITERATE

100.0% PROCESSED 10612 ITERATIONS 22 ANSWERS

SEARCH TIME: 00.00.01

L3 22 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
161.33
161.54

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FILE COVERS 1907 - 14 Sep 2005 VOL 143 ISS 12 FILE LAST UPDATED: 13 Sep 2005 (20050913/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13

L4 1 L3

=> d abs bib hitstr

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN

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GI
     For diagram(s), see printed CA Issue.
     The title compds. I and II [A = (substituted) 5- or 6-membered heteroarom.
AB
     ring or 5-, or 6-membered (hetero)cycloalkane ring; B =
     (hetero)cycloalkyl, (hetero)cycloalkenyl, or (hetero)aryl; L = alkylene;
     X, Y = a divalent linkage from the group consisting of alkylene,
     alkylene-OR3, alkylene-N(R3)COR4, alkylene-C(O)NR3R4, alkylene-N(R3)CO2R4,
     alkylene-N(R3)C(0)NR3R4, alkylene-C(0), O, C(0), N(R3), C(0)N(R3), S(0)k,
     and SO2N(R3); R1, R2 = H, alkyl, alkenyl, alkynyl, heteroalkyl, aryl,
     arylalkyl, NR6C(0)R5, C(0)R5, and NR5C(0)NR6; Rb = alkyl, aryl, OR7,
     C(0)R7, and C(0)NR7R8; R3, R4 = H, (hetero)alkyl, aryl, arylalkyl, etc.;
     R5, R6, R7, and R8 = H, alkyl, aryl, arylalkyl, etc.; p = 0-4; k = 0-2;
     with the proviso that X and Y are not both O, N(R3), S(O)k, or SO2N(R3)]
     were prepared as MCHR modulators in the treatment and/or prevention of
     endocrine, metabolic, cardiovascular, neurol., psychiatric,
     gastrointestinal, genitourinary and other disorders. For example,
     reaction of (2S)-2,3-dihydro-2-hydroxy-spiro[1H-indene-1,4'-piperidine]-1'-
     carboxylic acid 1,1-dimethylethyl ester (preparation given) with
     9-ethyl-9H-carbazole-3-carboxaldehyde yielded compound III. The prepared
     compds. were evaluated for MCHR2 modulation using a functional cell-based
     assay and most of them displayed potencies of 10 \mu M or less.
AN
     2004:290467 CAPLUS
     140:303531
DN
     Preparation of piperidinylmethyl carbazole derivatives as MCHR modulators
ΤI
     Chen, Xiaoqi; Dai, Kang; Fan, Pingchen; Fu, Ying; Li, Leping; Mihalic,
IN
     Jeffrey T.
     Tularik Inc., USA
PA
     PCT Int. Appl., 90 pp.
so
     CODEN: PIXXD2
DT
     Patent
     English
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                                                                         DATE
     PATENT NO.
                           KIND
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              GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
              LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM,
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                                                                         20030924
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PRAI US 2002-413335P
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     WO 2003-US30379
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os
     MARPAT 140:303531
     676606-93-2P 676607-01-5P 676607-06-0P
IT
     RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic
     preparation); THU (Therapeutic use); BIOL (Biological study); PREP
     (Preparation); RACT (Reactant or reagent); USES (Uses)
         (preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)
RN
     676606-93-2 CAPLUS
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CN Acetamide, N-[(3R)-1'-(9H-carbazol-3-ylmethyl)-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 676607-01-5 CAPLUS CN Spiro[1H-indene-1,4'-piperidine], 1'-(9H-carbazol-3-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-06-0 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-3-acetic acid, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-, monosodium salt (9CI) (CA INDEX NAME)

Na

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IT
     676606-89-6P 676606-90-9P 676606-92-1P
     676606-94-3P 676606-95-4P 676606-96-5P
     676606-97-6P 676606-98-7P 676606-99-8P
     676607-00-4P 676607-02-6P 676607-03-7P
     676607-04-8P 676607-05-9P 676607-07-1P
     676607-08-2P 676607-10-6P
     RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
        (preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)
RN
     676606-89-6 CAPLUS
     Spiro[1H-indene-1,4'-piperidin]-2-ol, 1'-[(9-ethyl-9H-carbazol-3-
CN
     yl)methyl]-2,3-dihydro-, (2S)- (9CI) (CA INDEX NAME)
```

Absolute stereochemistry.

RN 676606-90-9 CAPLUS CN Acetamide, N-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydro-2-hydroxyspiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-92-1 CAPLUS
CN Spiro[1H-indene-1,4'-piperidin]-3-ol, 3-ethyl-1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydro- (9CI) (CA INDEX NAME)

RN 676606-94-3 CAPLUS
CN Urea, N-ethyl-N'-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-95-4 CAPLUS

CN Carbamic acid, [1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]-, ethyl ester (9CI) (CA INDEX NAME)

RN 676606-96-5 CAPLUS

CN 4-Pyridineacetamide, N-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-97-6 CAPLUS

CN Spiro[1H-indene-1,4'-piperidin]-2-ol, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-, (2R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 676606-98-7 CAPLUS

CN 9H-Carbazole-9-carboxamide, 3-[(2,3-dihydro-2-hydroxyspiro[1H-indene-1,4'-piperidin]-1'-yl)methyl]-N,N-dimethyl- (9CI) (CA INDEX NAME)

RN 676606-99-8 CAPLUS

CN Spiro[1H-indene-1,4'-piperidin]-2-ol, 2,3-dihydro-1'-[(2,3,4,9-tetrahydro-1H-carbazol-6-yl)methyl]- (9CI) (CA INDEX NAME)

RN 676607-00-4 CAPLUS

CN 9H-Carbazole-9-carboxamide, 3-[[(3R)-3-(acetylamino)-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-1'-yl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN676607-02-6 CAPLUS

9H-Carbazole-9-carboxamide, 3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME) CN

RN

676607-03-7 CAPLUS
9H-Carbazole, 9-(1-oxopropyl)-3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME) CN

RN 676607-04-8 CAPLUS

CN 9H-Carbazole-9-carboxamide, N,N-dimethyl-3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-05-9 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine], 1'-[[9-(2-methoxyethyl)-9H-carbazol-3-yl]methyl]- (9CI) (CA INDEX NAME)

RN 676607-07-1 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-3-acetamide, 1'-(9H-carbazol-3-ylmethyl)2,3-dihydro-N-3-pyridinyl- (9CI) (CA INDEX NAME)

RN 676607-08-2 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-3-acetamide, 1'-(9H-carbazol-3-ylmethyl)2,3-dihydro-N-4-pyridinyl- (9CI) (CA INDEX NAME)

RN 676607-10-6 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-3-carboxamide, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-N-(1-methylethyl)- (9CI) (CA INDEX NAME)

IT 676607-27-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)

RN 676607-27-5 CAPLUS

CN Spiro[1H-indene-1,4'-piperidin]-3(2H)-one, 1'-[(9-ethyl-9H-carbazol-3-yl)methyl]- (9CI) (CA_INDEX_NAME)